

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG060724\
 Data File : BG061538.D
 Acq On : 7 Jun 2024 18:16
 Operator : MA/JU
 Sample : PB161251BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK251

Quant Time: Jun 07 22:53:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG052124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jun 06 12:15:45 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.871	152	25527	20.000	ng/u1	0.00
20) Naphthalene-d8	10.668	136	104543	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.516	164	82969	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.266	188	192389	20.000	ng/u1	0.00
79) Chrysene-d12	21.525	240	205223	20.000	ng/u1	0.00
88) Perylene-d12	24.634	264	233599	20.000	ng/u1	0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.329	96	2625	5.755	ng/uL	-0.03
4) Pyridine-d5	3.735	84	38406	23.667	ng/u1	-0.03
7) Phenol-d5	7.060	99	65195	31.100	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.201	67	37582	28.523	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.413	132	50449	32.518	ng/u1	0.00
15) 4-Methylphenol-d8	8.599	113	55943	31.297	ng/u1	0.00
21) Nitrobenzene-d5	9.028	128	27897	34.659	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.751	143	32827	34.848	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.309	165	64025	34.863	ng/u1	0.00
31) 4-Chloroaniline-d4	10.809	131	76743	31.941	ng/u1	0.00
46) Dimethylphthalate-d6	13.917	166	208925	32.467	ng/u1	0.00
49) Acenaphthylene-d8	14.205	160	226323	33.830	ng/u1	0.00
54) 4-Nitrophenol-d4	14.763	143	18997	23.879	ng/u1	0.00
60) Fluorene-d10	15.509	176	180418	32.475	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.656	200	24365	20.939	ng/u1	0.00
73) Anthracene-d10	17.366	188	295313	34.817	ng/u1	0.00
81) Pyrene-d10	19.663	212	349733	33.190	ng/u1	0.00
92) Benzo(a)pyrene-d12	24.422	264	391060	34.821	ng/u1	0.02

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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